

Welcome to STN International! Enter x:x

LOGINID:sssptal202txn

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

\* \* \* \* \* Welcome to STN International \* \* \* \* \*

NEWS	1		Web Page URLs for STN Seminar Schedule - N. America
NEWS	2		"Ask CAS" for self-help around the clock
NEWS	3	Feb 24	PCTGEN now available on STN
NEWS	4	Feb 24	TEMA now available on STN
NEWS	5	Feb 26	NTIS now allows simultaneous left and right truncation
NEWS	6	Feb 26	PCTFULL now contains images
NEWS	7	Mar 04	SDI PACKAGE for monthly delivery of multifile SDI results
NEWS	8	Mar 24	PATDPAFULL now available on STN
NEWS	9	Mar 24	Additional information for trade-named substances without structures available in REGISTRY
NEWS	10	Apr 11	Display formats in DGENE enhanced
NEWS	11	Apr 14	MEDLINE Reload
NEWS	12	Apr 17	Polymer searching in REGISTRY enhanced
NEWS	13	AUG 22	Indexing from 1927 to 1936 added to records in CA/CAPLUS
NEWS	14	Apr 21	New current-awareness alert (SDI) frequency in WPIDS/WPINDEX/WPIX
NEWS	15	Apr 28	RDISCLOSURE now available on STN
NEWS	16	May 05	Pharmacokinetic information and systematic chemical names added to PHAR
NEWS	17	May 15	MEDLINE file segment of TOXCENTER reloaded
NEWS	18	May 15	Supporter information for ENCOMPPAT and ENCOMPLIT updated
NEWS	19	May 19	Simultaneous left and right truncation added to WSCA
NEWS	20	May 19	RAPRA enhanced with new search field, simultaneous left and right truncation
NEWS	21	Jun 06	Simultaneous left and right truncation added to CBNB
NEWS	22	Jun 06	PASCAL enhanced with additional data
NEWS	23	Jun 20	2003 edition of the FSTA Thesaurus is now available
NEWS	24	Jun 25	HSDB has been reloaded
NEWS	25	Jul 16	Data from 1960-1976 added to RDISCLOSURE
NEWS	26	Jul 21	Identification of STN records implemented
NEWS	27	Jul 21	Polymer class term count added to REGISTRY
NEWS	28	Jul 22	INPADOC: Basic index (/BI) enhanced; Simultaneous Left and Right Truncation available
NEWS	29	AUG 05	New pricing for EUROPATFULL and PCTFULL effective August 1, 2003
NEWS	30	AUG 13	Field Availability (/FA) field enhanced in BEILSTEIN
NEWS	31	AUG 15	PATDPAFULL: one FREE connect hour, per account, in September 2003
NEWS	32	AUG 15	PCTGEN: one FREE connect hour, per account, in September 2003
NEWS	33	AUG 15	RDISCLOSURE: one FREE connect hour, per account, in September 2003
NEWS	34	AUG 15	TEMA: one FREE connect hour, per account, in September 2003
NEWS	35	AUG 18	Data available for download as a PDF in RDISCLOSURE
NEWS	36	AUG 18	Simultaneous left and right truncation added to PASCAL
NEWS	37	AUG 18	FROSTI and KOSMET enhanced with Simultaneous Left and Right Truncation
NEWS	38	AUG 18	Simultaneous left and right truncation added to ANABSTR

09/ 939,883

NEWS EXPRESS April 4 CURRENT WINDOWS VERSION IS V6.01a, CURRENT  
MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),  
AND CURRENT DISCOVER FILE IS DATED 01 APRIL 2003  
NEWS HOURS STN Operating Hours Plus Help Desk Availability  
NEWS INTER General Internet Information  
NEWS LOGIN Welcome Banner and News Items  
NEWS PHONE Direct Dial and Telecommunication Network Access to STN  
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that  
specific topic.

All use of STN is subject to the provisions of the STN Customer  
agreement. Please note that this agreement limits use to scientific  
research. Use for software development or design or implementation  
of commercial gateways or other similar uses is prohibited and may  
result in loss of user privileges and other penalties.

\* \* \* \* \* STN Columbus \* \* \* \* \*

FILE 'HOME' ENTERED AT 14:47:04 ON 28 AUG 2003 ✓

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 14:47:13 ON 28 AUG 2003

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file  
provided by InfoChem.

STRUCTURE FILE UPDATES: 26 AUG 2003 HIGHEST RN 573925-63-0

DICTIONARY FILE UPDATES: 26 AUG 2003 HIGHEST RN 573925-63-0

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2003

Please note that search-term pricing does apply when  
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP  
PROPERTIES for more information. See STNote 27, Searching Properties  
in the CAS Registry File, for complete details:

<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

=>

Uploading 09939883.str

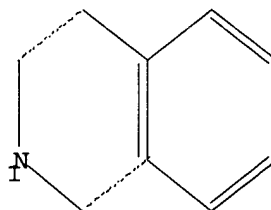
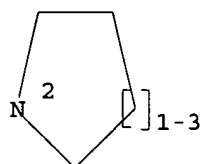
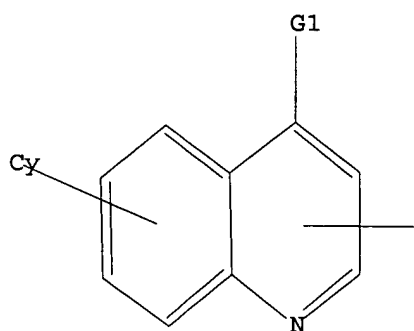
L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR

09/ 939,883



G1 [@1], [@2]

Structure attributes must be viewed using STN Express query preparation.

=> s l1 ful

FULL SEARCH INITIATED 14:47:37 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 11043 TO ITERATE

100.0% PROCESSED 11043 ITERATIONS

128 ANSWERS

SEARCH TIME: 00.00.01

L2 128 SEA SSS FUL L1

=> file caold

COST IN U.S. DOLLARS

SINCE FILE  
ENTRY

TOTAL  
SESSION

FULL ESTIMATED COST

148.15

148.36

FILE 'CAOLD' ENTERED AT 14:47:47 ON 28 AUG 2003

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

FILE COVERS 1907-1966

FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

09/ 939,883

This file supports REGISTRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> s l2

L3 0 L2

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.40

148.76

FILE 'CAPLUS' ENTERED AT 14:47:55 ON 28 AUG 2003

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 28 Aug 2003 VOL 139 ISS 9

FILE LAST UPDATED: 27 Aug 2003 (20030827/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l2

L4 2 L2

=> d l4 1- ibib abs hitstr

YOU HAVE REQUESTED DATA FROM 2 ANSWERS - CONTINUE? Y/(N):y

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2003 ACS on STN

ACCESSION NUMBER: 2002:185085 CAPLUS

DOCUMENT NUMBER: 136:247596

TITLE: Preparation of 7-aryl-4-(1-azacycloalkyl)quin(az)olines and analogs as NPY receptor antagonists

INVENTOR(S): Breu, Volker; Dautzenberg, Frank; Guerry, Philippe; Nettekoven, Matthias Heinrich; Pflieger, Philippe

PATENT ASSIGNEE(S): F. Hoffmann-La Roche A.-G., Switz.

SOURCE: PCT Int. Appl., 62 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2002020488	A2	20020314	WO 2001-EP10014	20010830
WO 2002020488	A3	20020516		

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,

GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,  
 LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PH, PL,  
 PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG,  
 UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM  
 RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,  
 DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF,  
 BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

US 2002052356 A1 20020502 US 2001-939883 20010827  
 AU 2002010474 A5 20020322 AU 2002-10474 20010830  
 BR 2001013710 A 20030603 BR 2001-13710 20010830  
 EP 1318981 A2 20030618 EP 2001-978324 20010830

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,  
 IE, SI, LT, LV, FI, RO, MK, CY, AL, TR

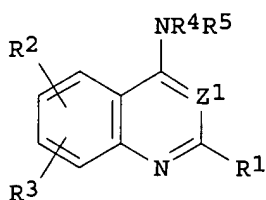
PRIORITY APPLN. INFO.:

EP 2000-119262 A 20000906  
 WO 2001-EP10014 W 20010830

OTHER SOURCE(S):

MARPAT 136:247596

GI



*Pregnant  
Version*

AB Title compds. [I; R1 = (cyclo)alkyl, CF3, aralkyl; R2 = H, halo, alkyl, alkoxy, etc.; R3 = (hetero)aryl; NR4R5 = (un)substituted heterocyclyl; Z1 = CH or N] were prepd. Thus, 4-chloro-7-iodo-2-methylquinoline was aminated by pyrrolidine and the product arylated by 3-ClC6H4B(OH)2 to give I [R1 = Me, R2 = H, R3 = 7-(3-chlorophenyl), R4R5 = (CH2)4, Z1 = CH].  
 Data for biol. activity of I were given.

IT 403849-18-3P 403849-19-4P 403849-20-7P  
 403849-21-8P 403849-22-9P 403849-23-0P  
 403849-24-1P 403849-25-2P 403849-26-3P  
 403849-27-4P 403849-28-5P 403849-29-6P  
 403849-30-9P 403849-31-0P 403849-32-1P  
 403849-33-2P 403849-34-3P 403849-35-4P  
 403849-43-4P 403849-44-5P 403849-45-6P  
 403849-46-7P 403849-48-9P 403849-49-0P  
 403849-90-1P 403849-92-3P 403849-96-7P  
 403849-98-9P 403850-00-0P 403850-02-2P  
 403850-04-4P 403850-06-6P 403850-08-8P  
 403850-10-2P 403850-12-4P 403850-14-6P  
 403850-16-8P 403850-17-9P 403850-18-0P  
 403850-19-1P 403850-20-4P 403850-22-6P  
 403850-23-7P 403850-25-9P 403850-27-1P  
 403850-29-3P 403850-31-7P 403850-32-8P  
 403850-33-9P 403850-35-1P 403850-37-3P  
 403850-39-5P 403850-40-8P 403850-41-9P  
 403850-43-1P 403850-45-3P 403850-47-5P  
 403850-50-0P 403850-52-2P 403850-54-4P  
 403850-56-6P 403850-58-8P 403850-60-2P  
 403850-62-4P 403850-64-6P 403850-65-7P  
 403853-09-8P 403853-11-2P

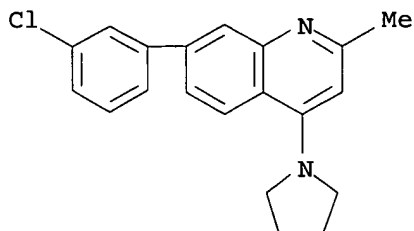
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of 7-aryl-4-(1-azacycloalkyl)quin(az)olines and analogs as NPY receptor antagonists)

09/ 939,883

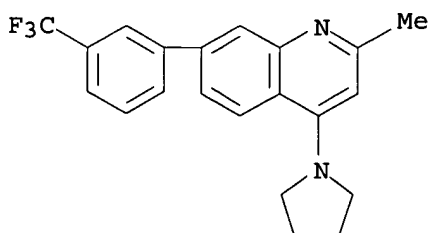
RN 403849-18-3 CAPLUS

CN Quinoline, 7-(3-chlorophenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



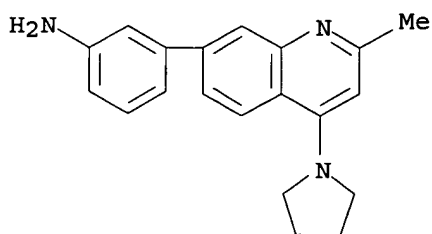
RN 403849-19-4 CAPLUS

CN Quinoline, 2-methyl-4-(1-pyrrolidinyl)-7-[3-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



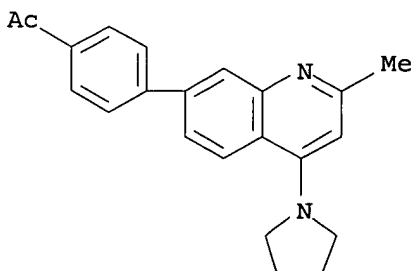
RN 403849-20-7 CAPLUS

CN Benzenamine, 3-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]- (9CI) (CA INDEX NAME)



RN 403849-21-8 CAPLUS

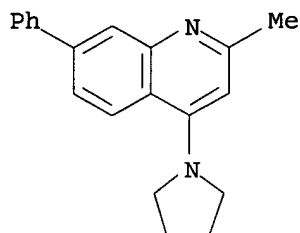
CN Ethanone, 1-[4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenyl]- (9CI) (CA INDEX NAME)



RN 403849-22-9 CAPLUS

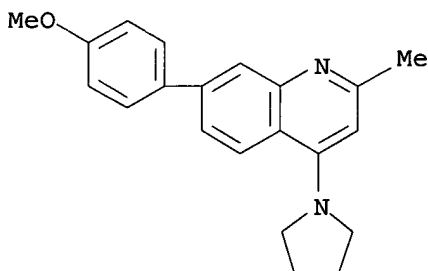
09/ 939,883

CN Quinoline, 2-methyl-7-phenyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



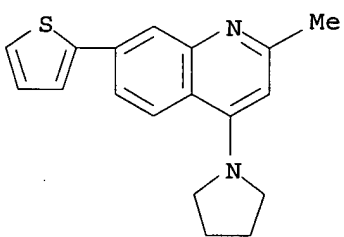
RN 403849-23-0 CAPLUS

CN Quinoline, 7-(4-methoxyphenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



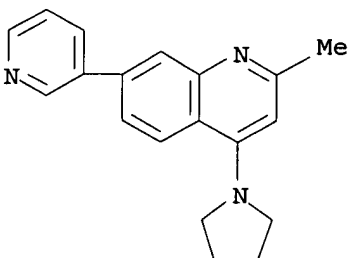
RN 403849-24-1 CAPLUS

CN Quinoline, 2-methyl-4-(1-pyrrolidinyl)-7-(2-thienyl)- (9CI) (CA INDEX NAME)



RN 403849-25-2 CAPLUS

CN Quinoline, 2-methyl-7-(3-pyridinyl)-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)

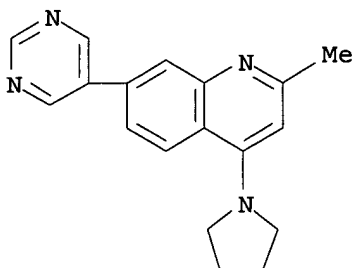


RN 403849-26-3 CAPLUS

CN Quinoline, 2-methyl-7-(5-pyrimidinyl)-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)

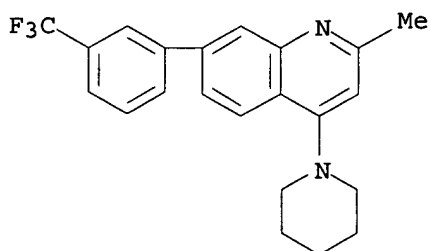
09/ 939,883

NAME)



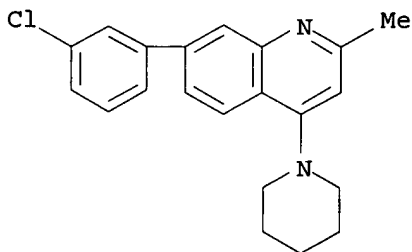
RN 403849-27-4 CAPLUS

CN Quinoline, 2-methyl-4-(1-piperidiny)-7-[3-(trifluoromethyl)phenyl] - (9CI)  
(CA INDEX NAME)



RN 403849-28-5 CAPLUS

CN Quinoline, 7-(3-chlorophenyl)-2-methyl-4-(1-piperidiny)- (9CI) (CA INDEX NAME)

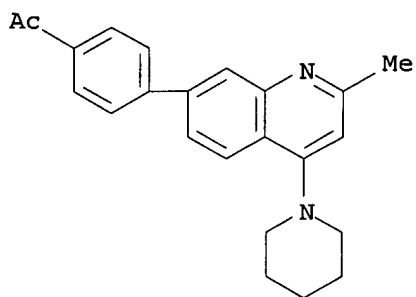


RN 403849-29-6 CAPLUS

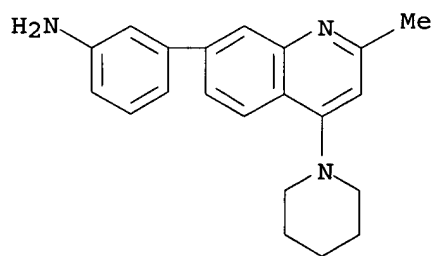
CN Ethanone, 1-[4-[2-methyl-4-(1-piperidiny)-7-quinolinyl]phenyl] - (9CI)  
(CA INDEX NAME)



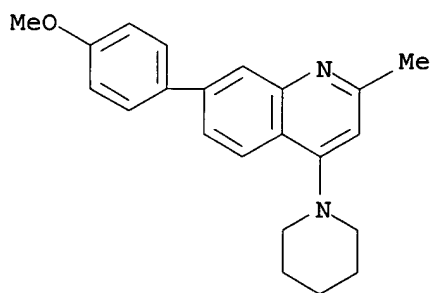
09/ 939,883



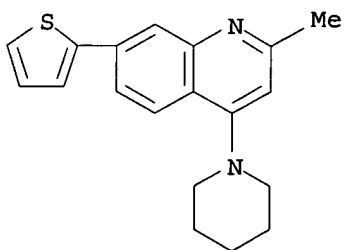
RN 403849-30-9 CAPLUS  
CN Benzenamine, 3-[2-methyl-4-(1-piperidinyl)-7-quinolinyl]- (9CI) (CA INDEX NAME)



RN 403849-31-0 CAPLUS  
CN Quinoline, 7-(4-methoxyphenyl)-2-methyl-4-(1-piperidinyl)- (9CI) (CA INDEX NAME)



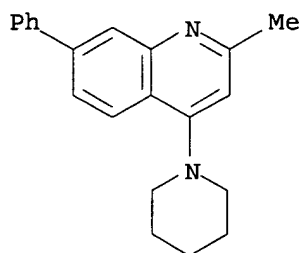
RN 403849-32-1 CAPLUS  
CN Quinoline, 2-methyl-4-(1-piperidinyl)-7-(2-thienyl)- (9CI) (CA INDEX NAME)



09/ 939,883

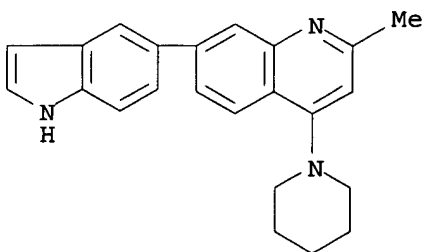
RN 403849-33-2 CAPLUS

CN Quinoline, 2-methyl-7-phenyl-4-(1-piperidinyl)- (9CI) (CA INDEX NAME)



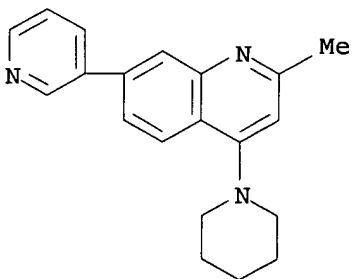
RN 403849-34-3 CAPLUS

CN Quinoline, 7-(1H-indol-5-yl)-2-methyl-4-(1-piperidinyl)- (9CI) (CA INDEX NAME)



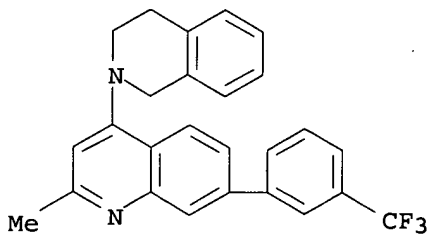
RN 403849-35-4 CAPLUS

CN Quinoline, 2-methyl-4-(1-piperidinyl)-7-(3-pyridinyl)- (9CI) (CA INDEX NAME)



RN 403849-43-4 CAPLUS

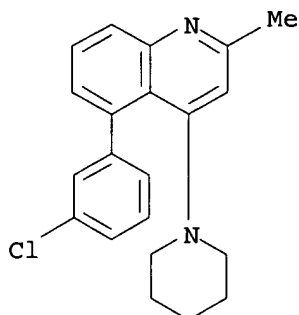
CN Quinoline, 4-(3,4-dihydro-2(1H)-isoquinolinyl)-2-methyl-7-[3-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



09/ 939,883

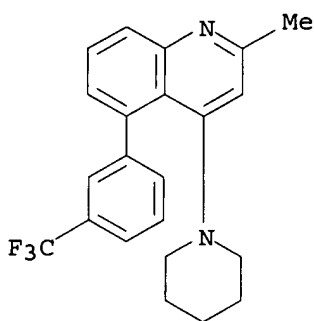
RN 403849-44-5 CAPLUS

CN Quinoline, 5-(3-chlorophenyl)-2-methyl-4-(1-piperidinyl)- (9CI) (CA INDEX NAME)



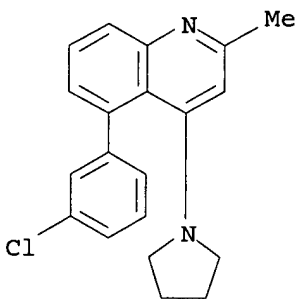
RN 403849-45-6 CAPLUS

CN Quinoline, 2-methyl-4-(1-piperidinyl)-5-[3-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)



RN 403849-46-7 CAPLUS

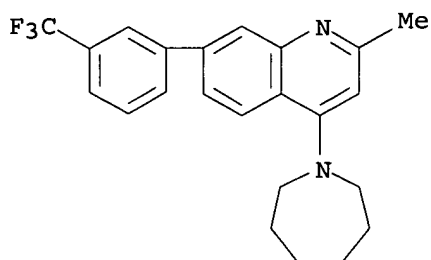
CN Quinoline, 5-(3-chlorophenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



RN 403849-48-9 CAPLUS

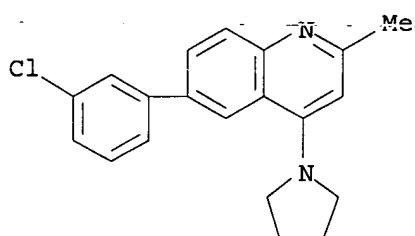
CN Quinoline, 4-(hexahydro-1H-azepin-1-yl)-2-methyl-7-[3-(trifluoromethyl)phenyl]- (9CI) (CA INDEX NAME)

09/ 939,883



RN 403849-49-0 CAPLUS

CN Quinoline, 6-(3-chlorophenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



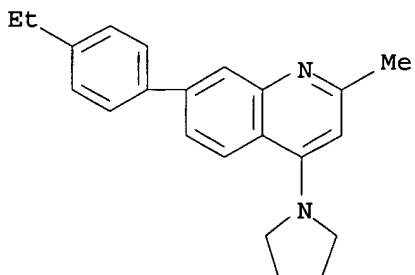
RN 403849-90-1 CAPLUS

CN Formic acid, compd. with 7-(4-ethylphenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403849-89-8

CMF C22 H24 N2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

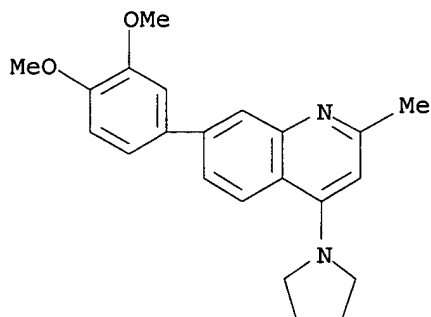
RN 403849-92-3 CAPLUS

CN Formic acid, compd. with 7-(3,4-dimethoxyphenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

09/ 939,883

CM 1

CRN 403849-91-2  
CMF C22 H24 N2 O2



CM 2

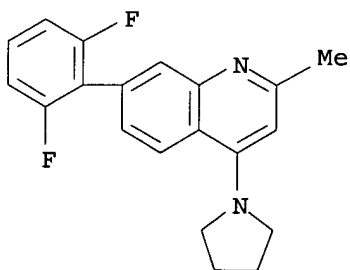
CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403849-96-7 CAPLUS  
CN Formic acid, compd. with 7-(2,6-difluorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403849-95-6  
CMF C20 H18 F2 N2



CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403849-98-9 CAPLUS  
CN Formic acid, compd. with 7-(2,4-dimethoxyphenyl)-2-methyl-4-(1-

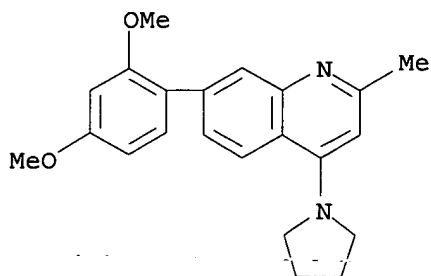
09/ 939,883

pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403849-97-8

CMF C22 H24 N2 O2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

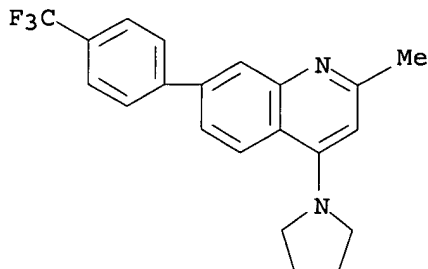
RN 403850-00-0 CAPLUS

CN Formic acid, compd. with 2-methyl-4-(1-pyrrolidinyl)-7-[4-(trifluoromethyl)phenyl]quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403849-99-0

CMF C21 H19 F3 N2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

RN 403850-02-2 CAPLUS

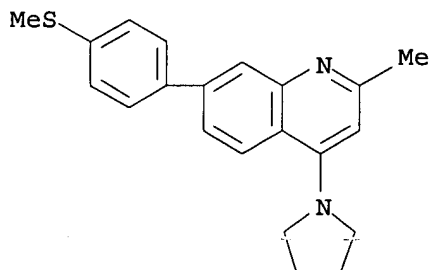
09/ 939,883

CN Formic acid, compd. with 2-methyl-7-[4-(methylthio)phenyl]-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-01-1

CMF C21 H22 N2 S



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

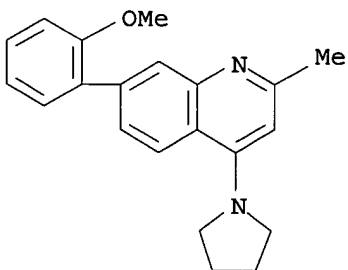
RN 403850-04-4 CAPLUS

CN Formic acid, compd. with 7-(2-methoxyphenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-03-3

CMF C21 H22 N2 O



CM 2

CRN 64-18-6

CMF C H2 O2

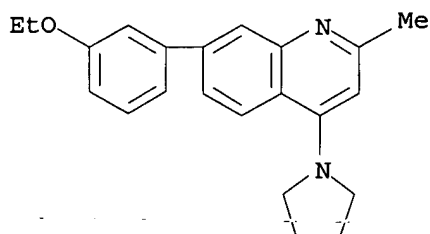
O=CH-OH

09/ 939,883

RN 403850-06-6 CAPLUS  
CN Formic acid, compd. with 7-(3-ethoxyphenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-05-5  
CMF C22 H24 N2 O



CM 2

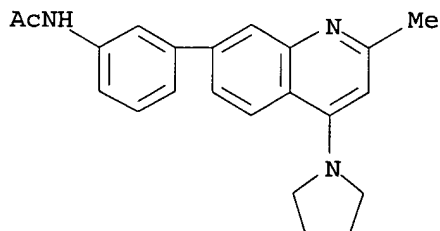
CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403850-08-8 CAPLUS  
CN Formic acid, compd. with N-[3-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenyl]acetamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-07-7  
CMF C22 H23 N3 O



CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403850-10-2 CAPLUS  
CN Formic acid, compd. with 2-methyl-4-(1-pyrrolidinyl)-7-[4-(trifluoromethoxy)phenyl]quinoline (1:1) (9CI) (CA INDEX NAME)

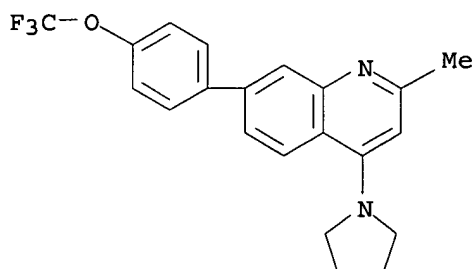


09/ 939,883

CM 1

CRN 403850-09-9

CMF C21 H19 F3 N2 O



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

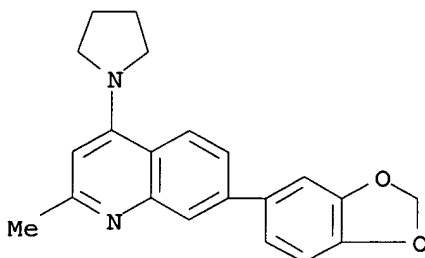
RN 403850-12-4 CAPLUS

CN Formic acid, compd. with 7-(1,3-benzodioxol-5-yl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-11-3

CMF C21 H20 N2 O2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

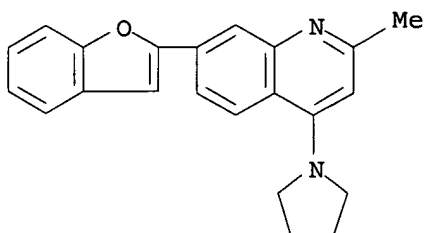
RN 403850-14-6 CAPLUS

CN Formic acid, compd. with 7-(2-benzofuranyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

09/ 939,883

CM 1

CRN 403850-13-5  
CMF C22 H20 N2 O



CM 2

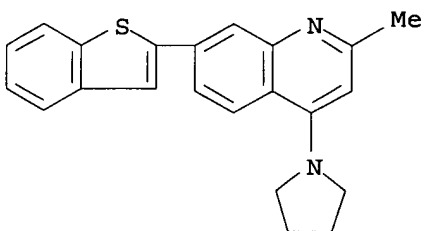
CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403850-16-8 CAPLUS  
CN Formic acid, compd. with 7-benzo[b]thien-2-yl-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-15-7  
CMF C22 H20 N2 S



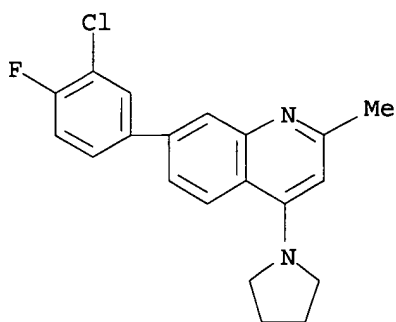
CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403850-17-9 CAPLUS  
CN Quinoline, 7-(3-chloro-4-fluorophenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI)  
(CA INDEX NAME)

09/ 939,883



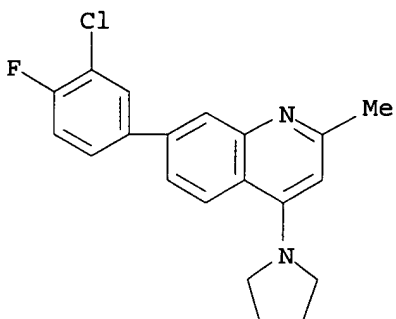
RN 403850-18-0 CAPLUS

CN Formic acid, compd. with 7-(3-chloro-4-fluorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-17-9

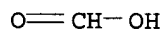
CMF C20 H18 Cl F N2



CM 2

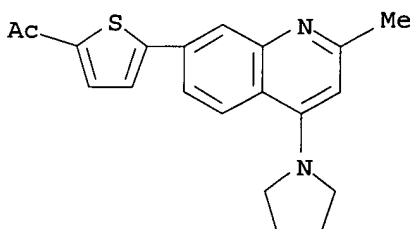
CRN 64-18-6

CMF C H2 O2



RN 403850-19-1 CAPLUS

CN Ethanone, 1-[5-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]-2-thienyl]- (9CI) (CA INDEX NAME)



RN 403850-20-4 CAPLUS

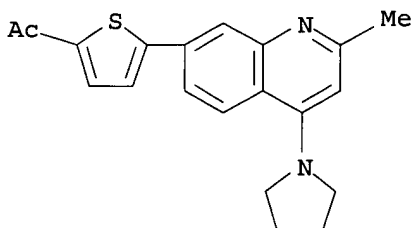
09/ 939,883

CN Formic acid, compd. with 1-[5-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]-2-thienyl]ethanone (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-19-1

CMF C20 H20 N2 O S



CM 2

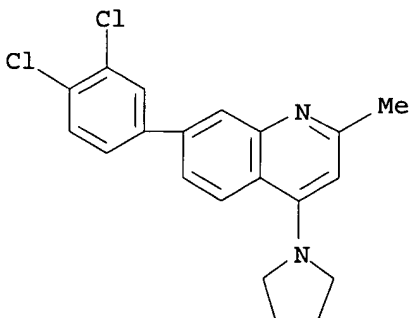
CRN 64-18-6

CMF C H2 O2

O=CH-OH

RN 403850-22-6 CAPLUS

CN Quinoline, 7-(3,4-dichlorophenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



RN 403850-23-7 CAPLUS

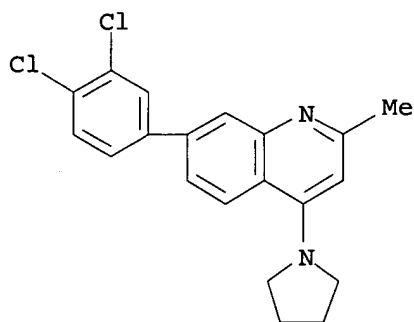
CN Formic acid, compd. with 7-(3,4-dichlorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-22-6

CMF C20 H18 Cl2 N2

09/ 939,883



CM 2

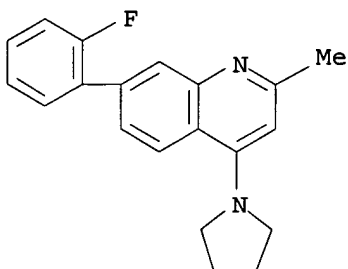
CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403850-25-9 CAPLUS  
CN Formic acid, compd. with 7-(2-fluorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-24-8  
CMF C20 H19 F N2



CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

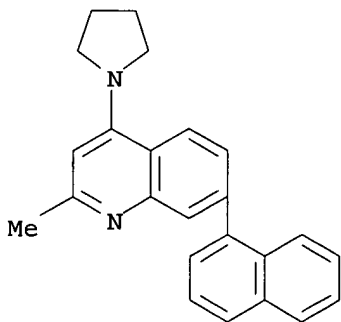
RN 403850-27-1 CAPLUS  
CN Formic acid, compd. with 2-methyl-7-(1-naphthalenyl)-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-26-0

09/ 939,883

CMF C24 H22 N2



CM 2

CRN 64-18-6  
CMF C H2 O2

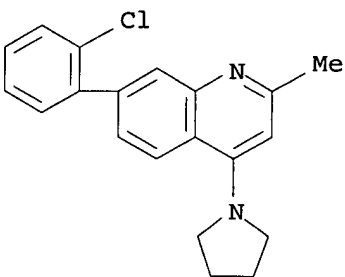
$$\text{O}=\text{CH}-\text{OH}$$

RN 403850-29-3 CAPLUS

CN      Formic acid, compd. with 7-(2-chlorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI)      (CA INDEX NAME)

CM 1

CRN 403850-28-2  
CMF C20 H19 Cl N2



CM 2

CRN 64-18-6  
CMF C H2 O2

$$\text{O}=\text{CH}-\text{OH}$$

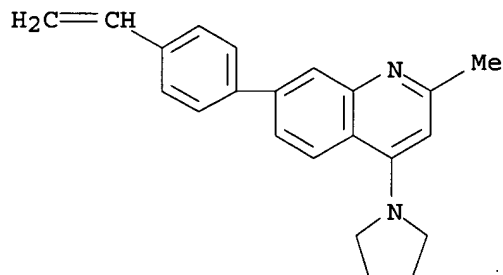
RN 403850-31-7 CAPLUS

CN    Formic acid, compd. with 7-(4-ethenylphenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI)    (CA INDEX NAME)

CM 1

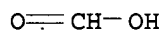
09/ 939,883

CRN 403850-30-6  
CMF C22 H22 N2

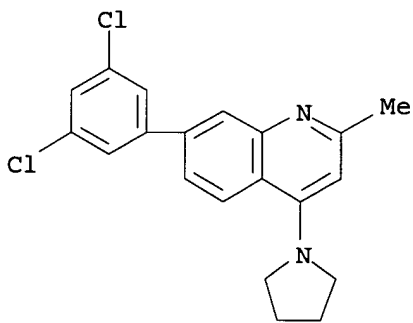


CM 2

CRN 64-18-6  
CMF C H2 O2



RN 403850-32-8 CAPLUS  
CN Quinoline, 7-(3,5-dichlorophenyl)-2-methyl-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)

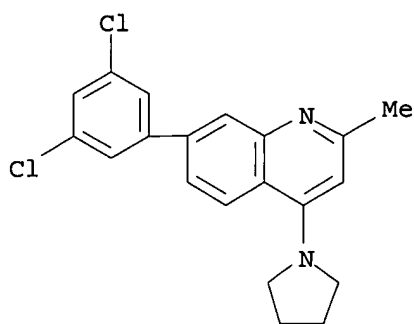


RN 403850-33-9 CAPLUS  
CN Formic acid, compd. with 7-(3,5-dichlorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-32-8  
CMF C20 H18 Cl2 N2

09/ 939,883



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

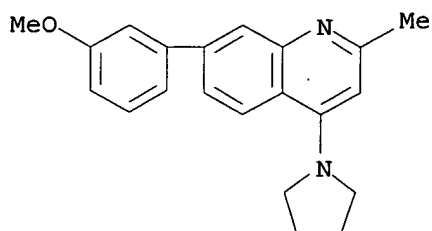
RN 403850-35-1 CAPLUS

CN Formic acid, compd. with 7-(3-methoxyphenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-34-0

CMF C21 H22 N2 O



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

RN 403850-37-3 CAPLUS

CN Benzoic acid, 3-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]-, ethyl ester, monoformate (9CI) (CA INDEX NAME)

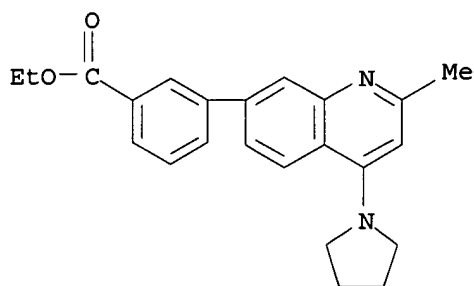
CM 1

CRN 403850-36-2

CMF C23 H24 N2 O2



09/ 939,883



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

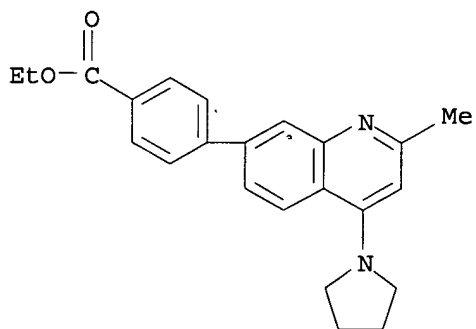
RN 403850-39-5 CAPLUS

CN Benzoic acid, 4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]-, ethyl ester, monoformate (9CI) (CA INDEX NAME)

CM 1

CRN 403850-38-4

CMF C23 H24 N2 O2



CM 2

CRN 64-18-6

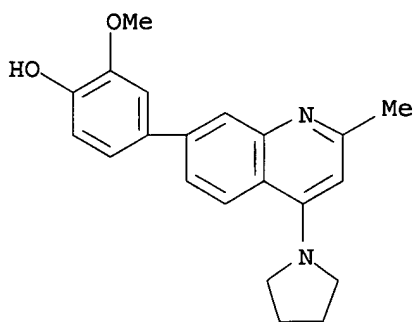
CMF C H2 O2

O=CH-OH

RN 403850-40-8 CAPLUS

CN Phenol, 2-methoxy-4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]- (9CI) (CA INDEX NAME)

09/ 939,883



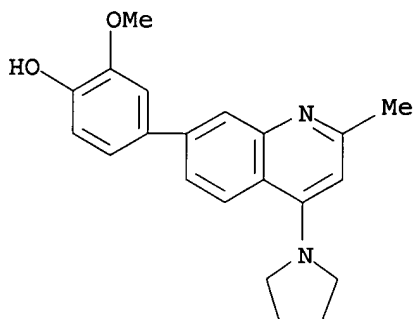
RN 403850-41-9 CAPLUS

CN Formic acid, compd. with 2-methoxy-4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-40-8

CMF C21 H22 N2 O2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

RN 403850-43-1 CAPLUS

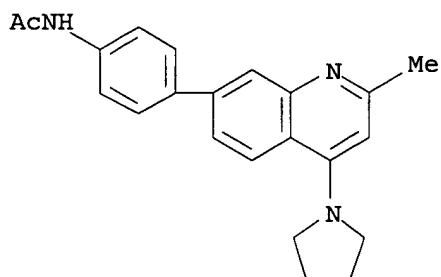
CN Formic acid, compd. with N-[4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenyl]acetamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-42-0

CMF C22 H23 N3 O

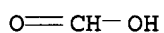
09/ 939,883



CM 2

CRN 64-18-6

CMF C H2 O2



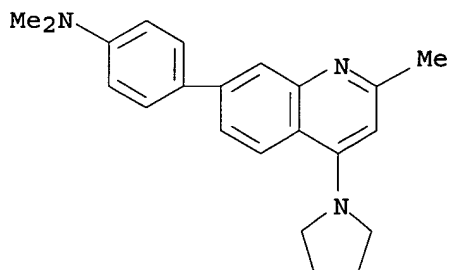
RN 403850-45-3 CAPLUS

CN Formic acid, compd. with N,N-dimethyl-4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]benzenamine (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-44-2

CMF C22 H25 N3



CM 2

CRN 64-18-6

CMF C H2 O2



RN 403850-47-5 CAPLUS

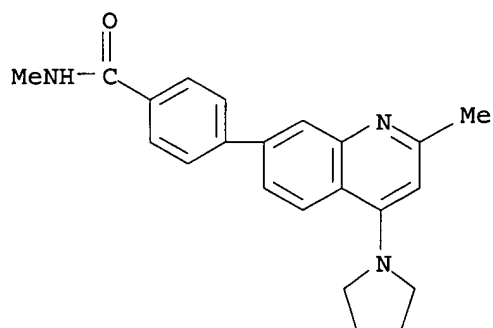
CN Formic acid, compd. with N-methyl-4-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]benzamide (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-46-4

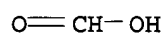
CMF C22 H23 N3 O

09/ 939,883



CM 2

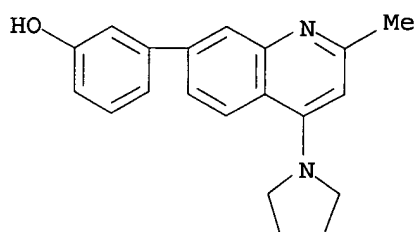
CRN 64-18-6  
CMF C H2 O2



RN 403850-50-0 CAPLUS  
CN Formic acid, compd. with 3-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenol (1:1) (9CI) (CA INDEX NAME)

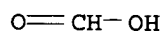
CM 1

CRN 403850-49-7  
CMF C20 H20 N2 O



CM 2

CRN 64-18-6  
CMF C H2 O2



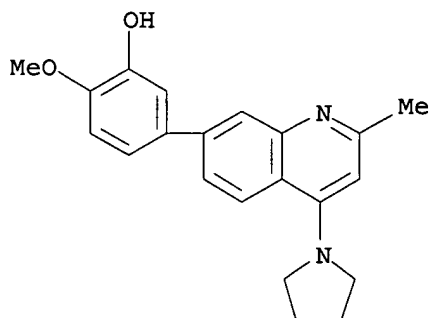
RN 403850-52-2 CAPLUS  
CN Formic acid, compd. with 2-methoxy-5-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenol (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-51-1

09/ 939,883

CMF C21 H22 N2 O2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

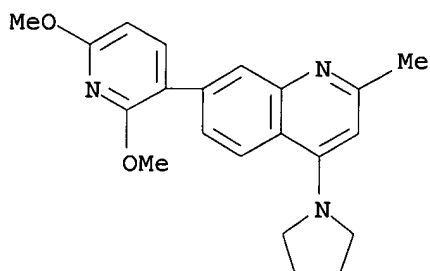
RN 403850-54-4 CAPLUS

CN Formic acid, compd. with 7-(2,6-dimethoxy-3-pyridinyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-53-3

CMF C21 H23 N3 O2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

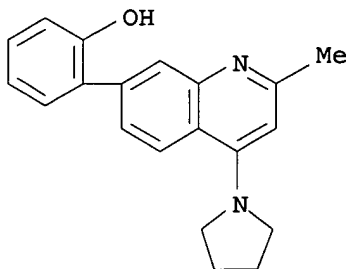
RN 403850-56-6 CAPLUS

CN Formic acid, compd. with 2-[2-methyl-4-(1-pyrrolidinyl)-7-quinolinyl]phenol (1:1) (9CI) (CA INDEX NAME)

CM 1

09/ 939,883

CRN 403850-55-5  
CMF C20 H20 N2 O



CM 2

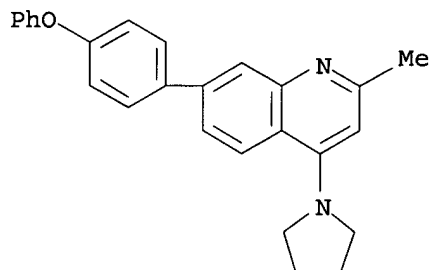
CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403850-58-8 CAPLUS  
CN Formic acid, compd. with 2-methyl-7-(4-phenoxyphenyl)-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-57-7  
CMF C26 H24 N2 O



CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

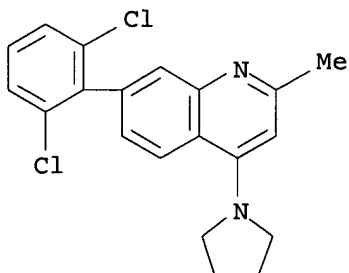
RN 403850-60-2 CAPLUS  
CN Formic acid, compd. with 7-(2,6-dichlorophenyl)-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

09/ 939,883

CM 1

CRN 403850-59-9

CMF C20 H18 Cl2 N2



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

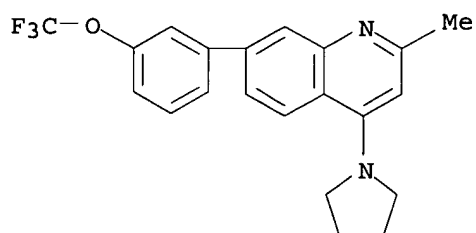
RN 403850-62-4 CAPLUS

CN Formic acid, compd. with 2-methyl-4-(1-pyrrolidinyl)-7-[3-(trifluoromethoxy)phenyl]quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-61-3

CMF C21 H19 F3 N2 O



CM 2

CRN 64-18-6

CMF C H2 O2

O=CH-OH

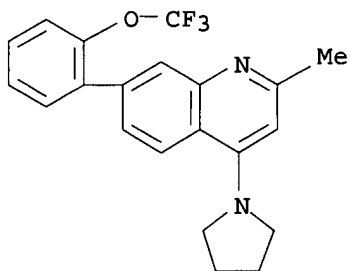
RN 403850-64-6 CAPLUS

CN Formic acid, compd. with 2-methyl-4-(1-pyrrolidinyl)-7-[2-(trifluoromethoxy)phenyl]quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

09/ 939,883

CRN 403850-63-5  
CMF C21 H19 F3 N2 O

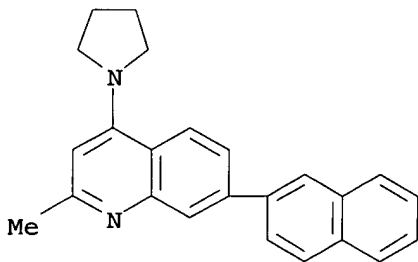


CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

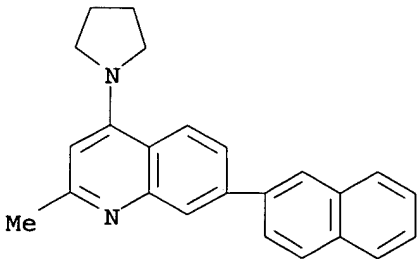
RN 403850-65-7 CAPLUS  
CN Quinoline, 2-methyl-7-(2-naphthalenyl)-4-(1-pyrrolidinyl)- (9CI) (CA INDEX NAME)



RN 403853-09-8 CAPLUS  
CN Formic acid, compd. with 2-methyl-7-(2-naphthalenyl)-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403850-65-7  
CMF C24 H22 N2





09/ 939,883

CM 2

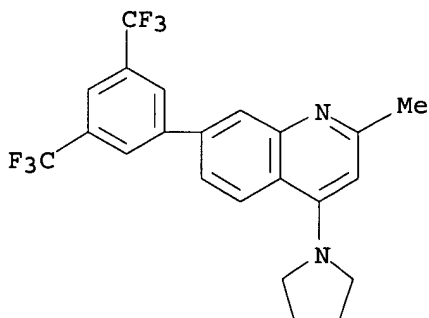
CRN 64-18-6  
CMF C H2 O2

O=CH-OH

RN 403853-11-2 CAPLUS  
CN Formic acid, compd. with 7-[3,5-bis(trifluoromethyl)phenyl]-2-methyl-4-(1-pyrrolidinyl)quinoline (1:1) (9CI) (CA INDEX NAME)

CM 1

CRN 403853-10-1  
CMF C22 H18 F6 N2



CM 2

CRN 64-18-6  
CMF C H2 O2

O=CH-OH

L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2003 ACS on STN  
ACCESSION NUMBER: 2002:31438 CAPLUS  
DOCUMENT NUMBER: 136:102370  
TITLE: Preparation of tetrahydropyridine or piperidine heterocyclic derivatives and their affinity for CRF receptors  
INVENTOR(S): Nakazato, Atsuro; Kumagai, Toshihito; Okubo, Taketoshi; Kameo, Kazuya  
PATENT ASSIGNEE(S): Taisho Pharmaceutical Co., Ltd., Japan  
SOURCE: PCT Int. Appl., 91 pp.  
CODEN: PIXXD2  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
-----	-----	-----	-----	-----

WO 2002002549 A1 20020110 WO 2001-JP5806 20010704  
 W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,  
 CO, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM,  
 HR, HU, ID, IL, IN, IS, JP, KE, KG, KR, KZ, LC, LK, LR, LS, LT,  
 LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU,  
 SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN,  
 YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM  
 RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,  
 DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF,  
 BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG  
 EP 1299378 A1 20030409 EP 2001-947819 20010704  
 R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,  
 IE, SI, LT, LV, FI, RO, MK, CY, AL, TR  
 NO 2002006125 A 20030204 NO 2002-6125 20021219  
 PRIORITY APPLN. INFO.: JP 2000-204021 A 20000705  
 JP 2000-270535 A 20000906  
 WO 2001-JP5806 W 20010704

OTHER SOURCE(S): MARPAT 136:102370

AB Tetrahydropyridine or piperidine heterocyclic derivs. with high affinity  
 for CRF receptors were prepd. E.g., 5-(4-carbamoyl-1,2,3,6-  
 tetrahydropyridin-1-yl)-2-(N-ethyl-2,4-dichloroanilino)-4-methylthiazole  
 was prepd. by bromination of 2-(N-ethyl-2,4-dichloroanilino)-4-  
 methylthiazole hydrochloride, followed by reaction with  
 5-carbamoyl-1,2,3,6-tetrahydropyridine hydrochloride.

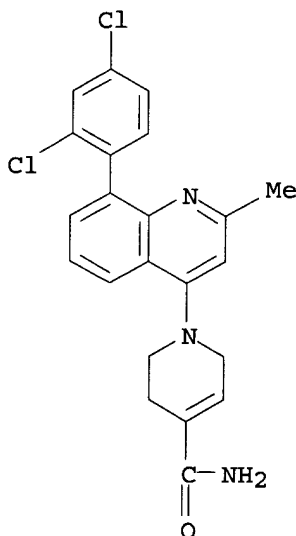
IT 388122-48-3P 388122-49-4P 388122-50-7P  
 388122-52-9P 388122-53-0P 388122-54-1P  
 388122-55-2P 388122-56-3P 388122-57-4P  
 388122-58-5P 388122-59-6P 388122-60-9P  
 388122-61-0P 388122-62-1P 388122-63-2P  
 388122-64-3P 388122-65-4P 388122-67-6P  
 388122-68-7P 388122-69-8P 388122-70-1P  
 388122-71-2P 388122-72-3P 388122-73-4P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU  
 (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES  
 (Uses)

(prepn. of tetrahydropyridine or piperidine heterocyclic derivs. and  
 their affinity for CRF receptors)

RN 388122-48-3 CAPLUS

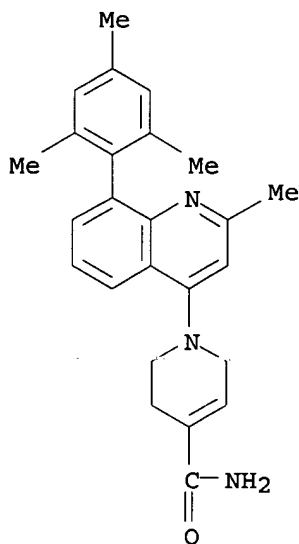
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-  
 1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



09/ 939,883

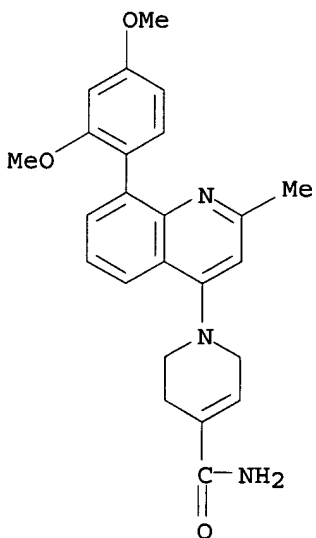
RN 388122-49-4 CAPLUS

CN 4-Pyridinecarboxamide, 1,2,3,6-tetrahydro-1-[2-methyl-8-(2,4,6-trimethylphenyl)-4-quinolinyl]- (9CI) (CA INDEX NAME)



RN 388122-50-7 CAPLUS

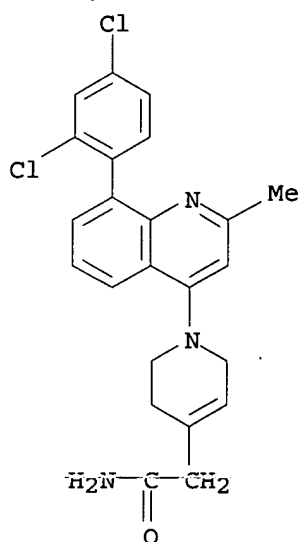
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dimethoxyphenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-52-9 CAPLUS

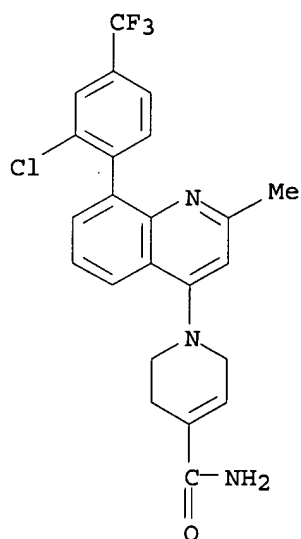
CN 4-Pyridineacetamide, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-53-0 CAPLUS

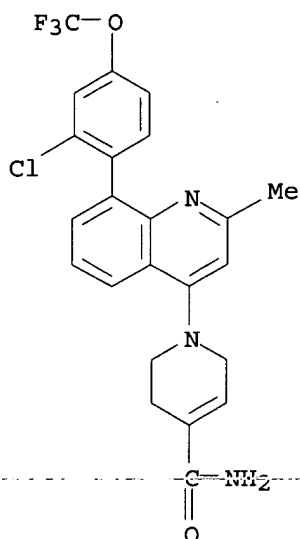
CN 4-Pyridinecarboxamide, 1-[8-[2-chloro-4-(trifluoromethyl)phenyl]-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-54-1 CAPLUS

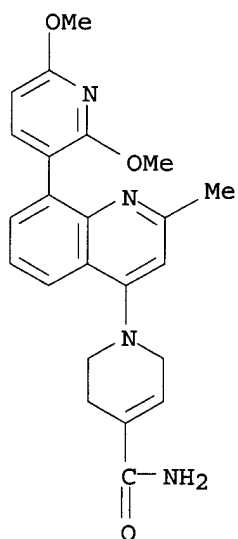
CN 4-Pyridinecarboxamide, 1-[8-[2-chloro-4-(trifluoromethoxy)phenyl]-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-55-2 CAPLUS

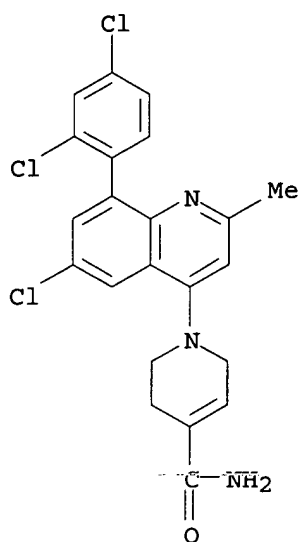
CN 4-Pyridinecarboxamide, 1-[8-(2,6-dimethoxy-3-pyridinyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-56-3 CAPLUS

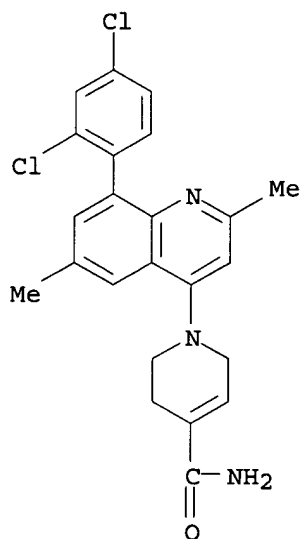
CN 4-Pyridinecarboxamide, 1-[6-chloro-8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-57-4 CAPLUS

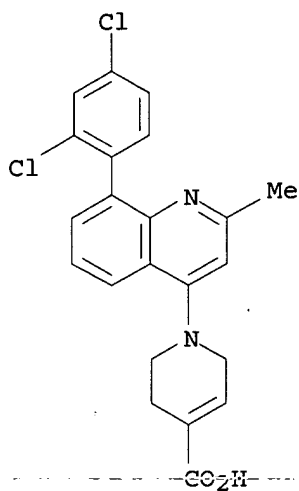
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-2,6-dimethyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-58-5 CAPLUS

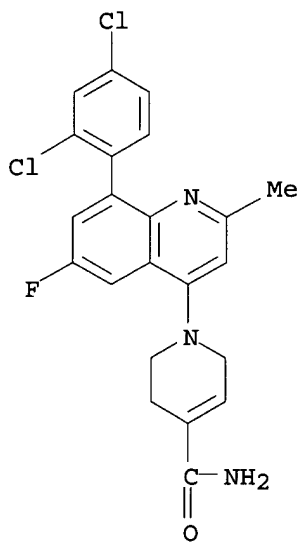
CN 4-Pyridinecarboxylic acid, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-59-6 CAPLUS

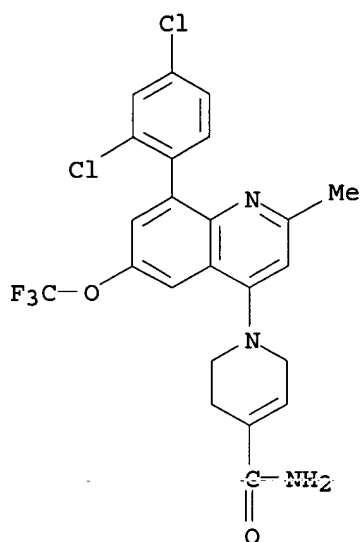
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-6-fluoro-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-60-9 CAPLUS

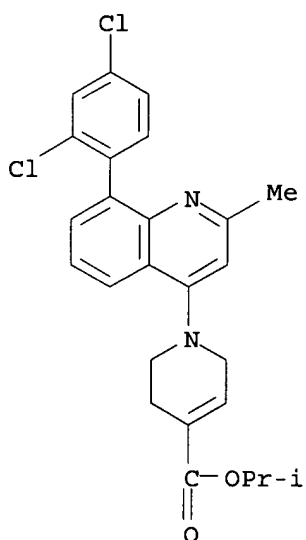
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-2-methyl-6-(trifluoromethoxy)-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-61-0 CAPLUS

CN 4-Pyridinecarboxylic acid, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro-, 1-methylethyl ester (9CI) (CA INDEX NAME)

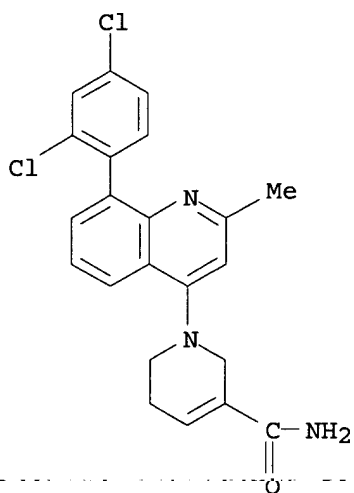


RN 388122-62-1 CAPLUS

CN 3-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,5,6-tetrahydro- (9CI) (CA INDEX NAME)

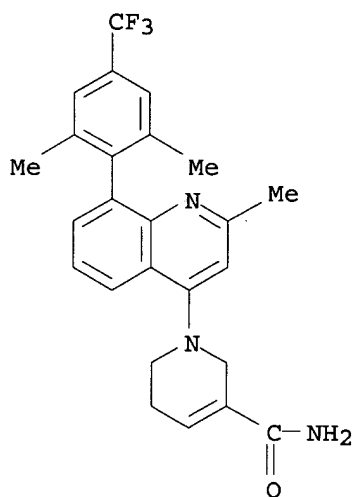


09/ 939,883



RN 388122-63-2 CAPLUS

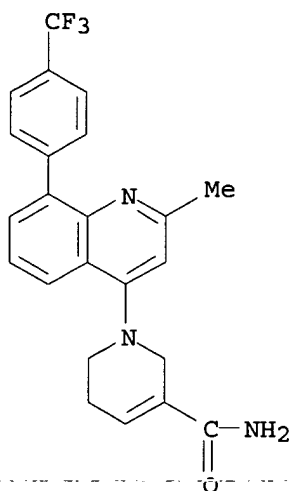
CN 3-Pyridinecarboxamide, 1-[8-[2,6-dimethyl-4-(trifluoromethyl)phenyl]-2-methyl-4-quinolinyl]-1,2,5,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-64-3 CAPLUS

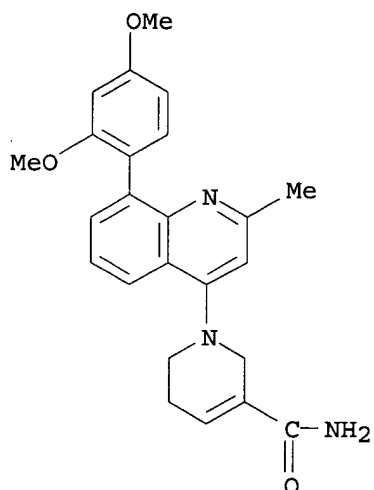
CN 3-Pyridinecarboxamide, 1,2,5,6-tetrahydro-1-[2-methyl-8-[4-(trifluoromethyl)phenyl]-4-quinolinyl]- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-65-4 CAPLUS

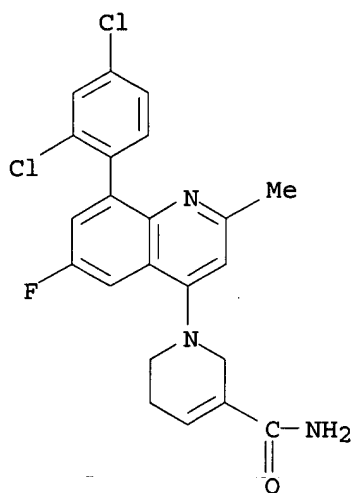
CN 3-Pyridinecarboxamide, 1-[8-(2,4-dimethoxyphenyl)-2-methyl-4-quinolinyl]-1,2,5,6-tetrahydro- (9CI) (CA INDEX NAME)



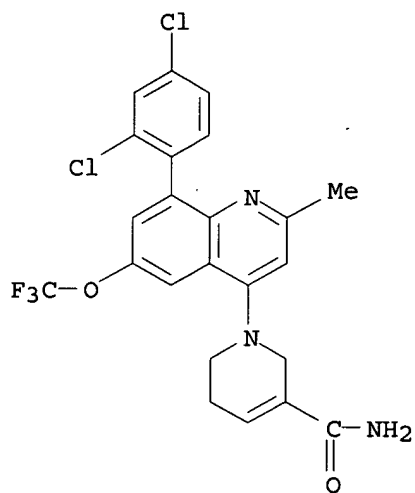
RN 388122-67-6 CAPLUS

CN 3-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-6-fluoro-2-methyl-4-quinolinyl]-1,2,5,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



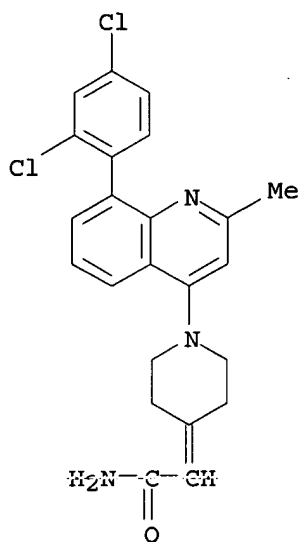
RN 388122-68-7 CAPLUS  
CN 3-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-2-methyl-6-(trifluoromethoxy)-4-quinolinyl]-1,2,5,6-tetrahydro-, monohydrochloride (9CI) (CA INDEX NAME)



● HCl

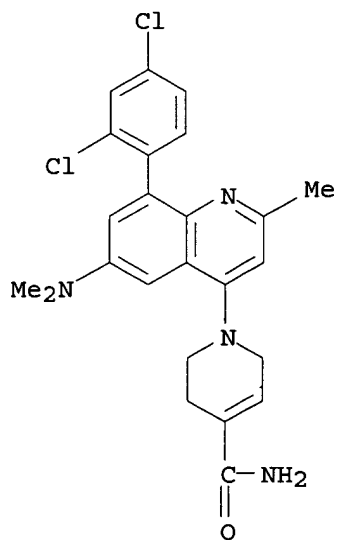
RN 388122-69-8 CAPLUS  
CN Acetamide, 2-[1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-4-piperidinylidene]- (9CI) (CA INDEX NAME)

09/ 939,883



RN 388122-70-1 CAPLUS

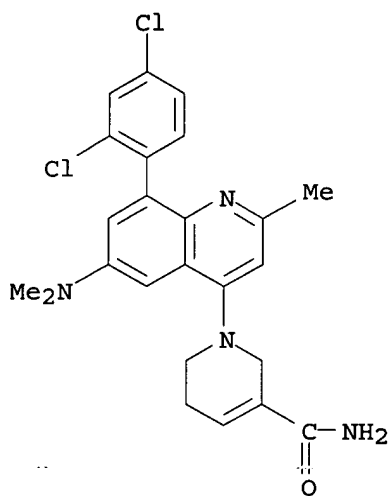
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-6-(dimethylamino)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-71-2 CAPLUS

CN 3-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-6-(dimethylamino)-2-methyl-4-quinolinyl]-1,2,5,6-tetrahydro-, monohydrochloride (9CI) (CA INDEX NAME)

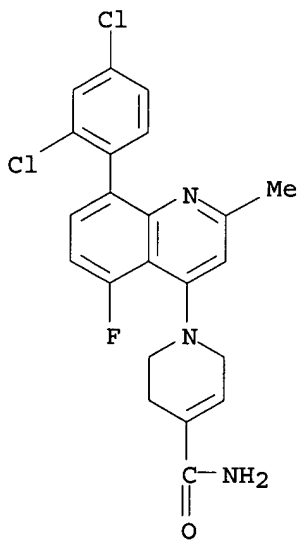
09/ 939,883



● HCl

RN 388122-72-3 CAPLUS

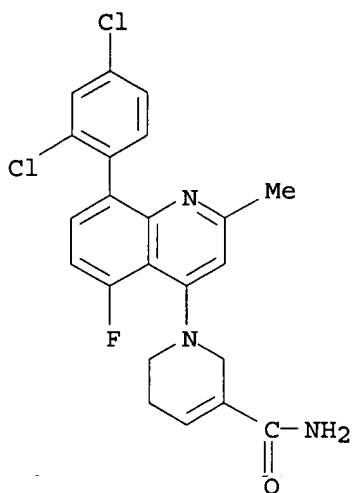
CN 4-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-5-fluoro-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388122-73-4 CAPLUS

CN 3-Pyridinecarboxamide, 1-[8-(2,4-dichlorophenyl)-5-fluoro-2-methyl-4-quinolinyl]-1,2,5,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



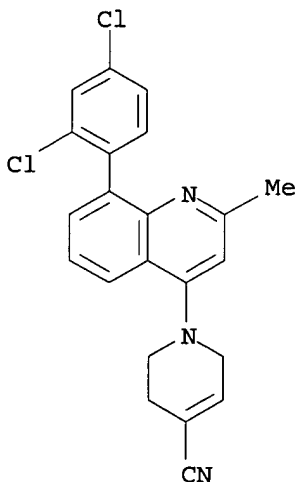
IT 388123-52-2P 388123-59-9P 388123-60-2P  
388123-61-3P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(prepn. of tetrahydropyridine or piperidine heterocyclic derivs. and their affinity for CRF receptors)

RN 388123-52-2 CAPLUS

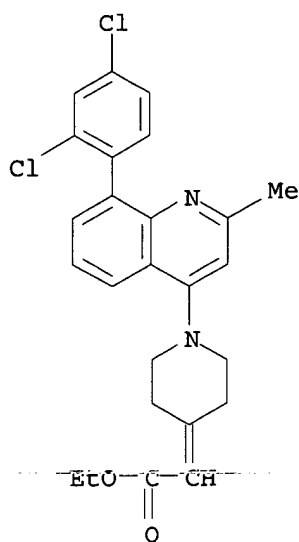
CN 4-Pyridinecarbonitrile, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)



RN 388123-59-9 CAPLUS

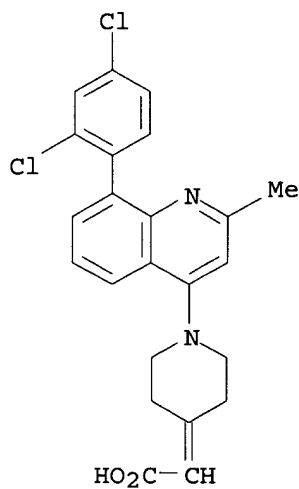
CN Acetic acid, [1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-4-piperidinylidene]-, ethyl ester (9CI) (CA INDEX NAME)

09/ 939,883



RN 388123-60-2 CAPLUS

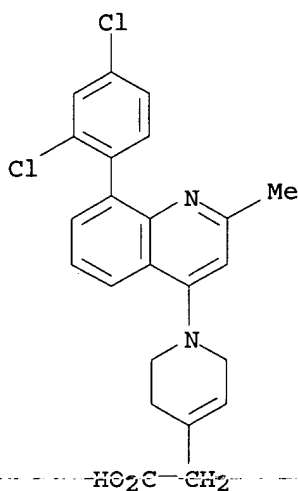
CN Acetic acid, [1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-4-piperidinylidene]- (9CI) (CA INDEX NAME)



RN 388123-61-3 CAPLUS

CN 4-Pyridineacetic acid, 1-[8-(2,4-dichlorophenyl)-2-methyl-4-quinolinyl]-1,2,3,6-tetrahydro- (9CI) (CA INDEX NAME)

09/ 939,883



REFERENCE COUNT: 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> d his

(FILE 'HOME' ENTERED AT 14:47:04 ON 28 AUG 2003)

FILE 'REGISTRY' ENTERED AT 14:47:13 ON 28 AUG 2003

L1 STRUCTURE UPLOADED  
L2 128 S L1 FUL

FILE 'CAOLD' ENTERED AT 14:47:47 ON 28 AUG 2003

L3 0 S L2

FILE 'CAPLUS' ENTERED AT 14:47:55 ON 28 AUG 2003

L4 2 S L2

=> log y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

9.91

158.67

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-1.30

-1.30

STN INTERNATIONAL LOGOFF AT 14:48:56 ON 28 AUG 2003